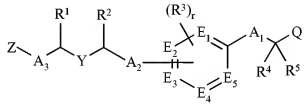
Amendments to the Claims

IN THE CLAIMS

Please cancel Claims 4, 8, 9, 32, 45, 46, and 48.

Amendments to the Claims

1. (Original). A compound having a formula I,



I

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 is: a bond, CH_2, O or S, and wherein A_1 and R^4 or A_1 and R^5 together being a 3- to 6-membered carbocyclyl when A_1 is a carbon;

A_2 and A_3 are independently: CH_2, O or S;

E_1, E_2, E_3, E_4 and E_5 are each CH or substituted carbon bearing A_2 and R^3; or at least one of E_1, E_2, E_3, E_4 and E_5 is nitrogen and each of others being CH or substituted carbon bearing A_2 and R^3;

Q is: $-C(O)OR^6$, or R^{6A};

Y is: a bond, C_1-C_6 alkyl or C_3-C_6 cycloalkyl;

Z is: a) aryl;

b) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,

c) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or

d) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R^7;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: 1, 2, 3, or 4;

R^1 and R^2 are each independently:

hydrogen,

haloalkyl,

C_1-C_6 alkyl,

$(CH_2)_nC_3$-C_8 cycloalkyl, or

R^1 and R^2 form a 4- to 8-membered nonaromatic carbocyclic ring; and

wherein at least one of R^1 and R^2 is alkyl or cycloalkyl, and;

R^3 is: hydrogen,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

C_1-C_6 alkyl,

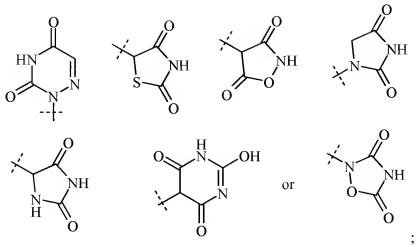
C_1-C_6 alkoxy, or

C_3-C_8 cycloalkyl;

R^4 and R^5 are each independently: hydrogen or C_1-C_6 alkyl;

R^6 is: hydrogen, C_1-C_6 alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,



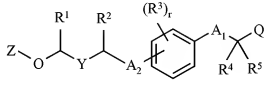
;

R^7 is: hydrogen,
oxo,
nitro,
cyano,
hydroxyl,
halo,
haloalkyl,
haloalkyloxy,
aryloxy,
arylalkyl,
aminoalkyl,
C_1-C_6 alkyl,
C_1-C_6 alkoxy,
$(CH_2)_nC_3$-C_8 cycloalkyl,
$C(O)R^9$,
$C(O)OR^9$,
$C(=NOR^8)R^9$,
$CR^8(OH)R^9$,
$C[=C(R^8)_2]R^9$,
OR^9,
SR^9 or
$S(O)_nR^9$;

R^8 is: hydrogen or C_1-C_6 alkyl; and

R^9 is: hydrogen,
C_1-C_6 alkyl,
C_3-C_8 cycloalkyl,
aryl,
heteroaryl or
heterocyclyl,
wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:
hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy and C_3-C_8 cycloalkyl.

2. (Previously Presented). The compound of Claim 1, wherein the compound is represented by a compound of formula II,



II

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 is: a bond, CH_2, O or S, and wherein A_1 and R^4 or A_1 and R^5 together being a 3- to 6-membered carbocyclyl when A_1 is a carbon;

A_2 is: O or S or CH_2;

Q is: $-C(O)OR^6$, or R^{6A};

Y is: a bond, C_1-C_6 alkyl or C_3-C_6 cycloalkyl;

Z is: a) aryl;
b) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,
c) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or

d) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and
wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R^7;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: 1, 2, 3, or 4;

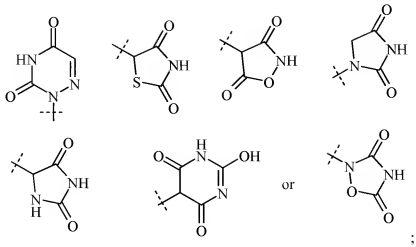
R^1 and R^2 are each independently:
hydrogen,
haloalkyl,
C_1-C_6 alkyl,
$(CH_2)_nC_3$-C_8 cycloalkyl, or
R^1 and R^2 form a 4- to 8-membered nonaromatic carbocyclic ring; and
wherein at least one of R^1 and R^2 is alkyl or cycloalkyl, and;

R^3 is: hydrogen,
nitro,
cyano,
hydroxyl,
halo,
haloalkyl,
haloalkyloxy,
aryloxy,
C_1-C_6 alkyl,
C_1-C_6 alkoxy, or
C_3-C_8 cycloalkyl;

R^4 and R^5 are each independently: hydrogen or C_1-C_6 alkyl;

R^6 is: hydrogen, C_1-C_6 alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,



;

R^7 is: hydrogen,
oxo,
nitro,
cyano,
hydroxyl,
halo,
haloalkyl,
haloalkyloxy,
aryloxy,
arylalkyl,
aminoalkyl,
C_1-C_6 alkyl,
C_1-C_6 alkoxy,
$(CH_2)_nC_3$-C_8 cycloalkyl,
$C(O)R^9$,
$C(O)OR^9$,
$C(=NOR^8)R^9$,
$CR^8(OH)R^9$,
$C[=C(R^8)_2]R^9$,
OR^9,
SR^9 or
$S(O)_nR^9$;

R^8 is: hydrogen or C_1-C_6 alkyl; and

R^9 is: hydrogen,

C_1-C_6 alkyl,

C_3-C_8 cycloalkyl,

aryl,

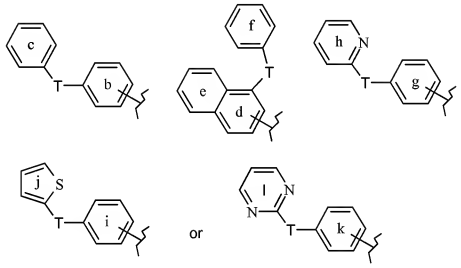
heteroaryl or

heterocyclyl,

wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy and C_3-C_8 cycloalkyl.

3. (Original). The compound of Claim 2, wherein Z is optionally substituted phenyl or naphthyl, furanyl, imidazolyl, indolyl, oxazolyl, isoxazolyl, pyridyl, pyrrolyl, thiazolyl, thiophenyl, benzofuranyl, benzothiophenyl, benzoisoxazolyl, quinolinyl, isoquinolinyl or a structural formula selected from following:



wherein T is:

a bond, $-(CH_2)_qO-$, $-O(CH_2)_q-$, $-C(O)(CH_2)_q-$, $-(CH_2)_qC(O)-$, $-(CH_2)_qS-$, $-S(CH_2)_q-$, $S[O]_p$, $-(C_1$-C_3 alkyl$)-$, $-(CH_2)_qC(=CH_2)-$, $-C(=CH_2)(CH_2)_q-$, $-(CH_2)_qC(=NOH)-$, $-C(=NOH)(CH_2)_q-$, $-(CH_2)_qC(=NOCH_3)-$, $-C(=NOCH_3)(CH_2)_q-$, $-CH(OH)(CH_2)_q-$, or $-(CH_2)_qCH(OH)-$,

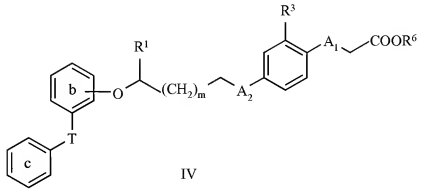
q is: 0, 1, 2 or 3; and

rings b to l are each optionally substituted with one or more groups independently selected from the group consisting of:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1-C_6 alkyl, C_1-C_6 alkoxy and $(CH_2)_nC_3$-C_8 cycloalkyl.

4. (Canceled)

5. (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula IV,



IV

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 and A_2 are respectively:

O and O,

CH_2 and O,

CH_2 and S,

O and S or

S and O;

m is: 1 or 2;

R^1 is: C_1-C_3 alkyl;

R^3 is: hydrogen, halo or C_1-C_6 alkyl;

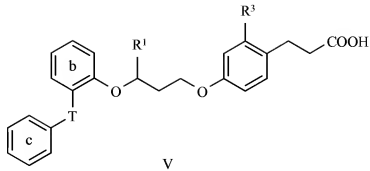
R^6 and R^9 are each independently: hydrogen or C_1-C_6 alkyl;

T is: a bond, -O-, -C(O)-, -S(O) –$S(O)_2$-, -C(=CH_2)-, -C(=NOH)- or -CH(OH)-; and

rings b and c are each optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1-C_6 alkyl, C_1-C_6 alkoxy and $(CH_2)_nC_3$-C_8 cycloalkyl.

6. (Previously Presented). The compound of Claim 5, wherein the compound is represented by structural formula V,



V

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

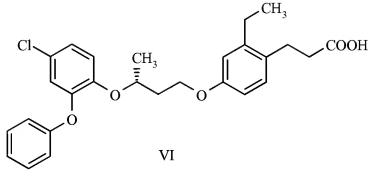
T is: a bond, -O- or -C(O)-;

R^1 is: methyl, ethyl or cyclopropyl;

R^3 is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF_3, OCF_3, methyl, ethyl, isopropyl, $N(CH_3)_2$, $S(O)_2CH_3$, methoxy and cyclopropyl.

7. (Original). The compound of Claim 6, wherein the compound is represented by a structural formula VI,



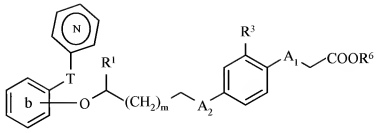
VI

or a pharmaceutically acceptable salt, solvate or hydrate thereof.

8. (Canceled)

9. (Canceled)

10. (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula VIII,



VIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 and A_2 are respectively:

O and O,

CH_2 and O,

CH_2 and S,

O and S or

S and O;

m is: 1 or 2;

R^1 is: C_1-C_3 alkyl; and

R^3 is: hydrogen, halo or C_1-C_6 alkyl;

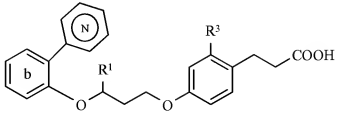
R^6 and R^9 are each independently: hydrogen or C_1-C_6 alkyl;

T is: a bond, -O-, -C(O)-, -S(O) –$S(O)_2$-, -C(=CH_2)-, -C(=NOH)- or -CH(OH)-; and

ring b is optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1-C_6 alkyl, C_1-C_6 alkoxy and $(CH_2)_nC_3$-C_8 cycloalkyl.

11. (Previously Presented). The compound of Claim 10, wherein the compound is represented by structural formula IX,



IX

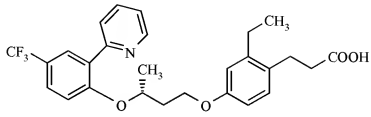
or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

R^1 is C_1-C_3 alkyl;

R^3 is: hydrogen, halo or C_1-C_4 alkyl;

ring b is optionally substituted with one or more groups independently selected from the group consisting of: hydrogen, halo, haloalkyl, haloalkyloxy and C_1-C_6 alkyl.

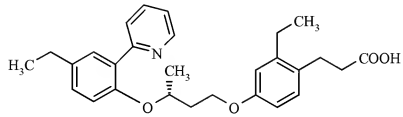
12. (Previously Presented). The compound of Claim 11, wherein the compound is represented by structural formula X,



X

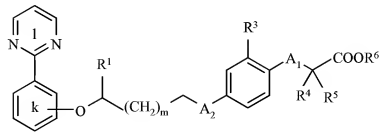
or a pharmaceutically acceptable salt, solvate or hydrate thereof.

13. (Previously Presented). The compound of Claim 11, wherein the compound is represented by structural formula XI,



XI

14. (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula XII,



XII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:
A_1 and A_2 are respectively:

O and O,

CH_2 and O,

CH_2 and S,

O and S or

S and O;

m is: 1 or 2;

R^1 is: C_1-C_3 alkyl; and

R^3 is: hydrogen, halo or C_1-C_6 alkyl;

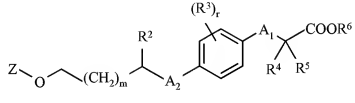
R^4, R^5, R^6 and R^9 are each independently: hydrogen or C_1-C_6 alkyl;

rings k and l are each optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1-C_6 alkyl, C_1-C_6 alkoxy and $(CH_2)_nC_3$-C_8 cycloalkyl.

15. (Canceled)

16. (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula XIII,

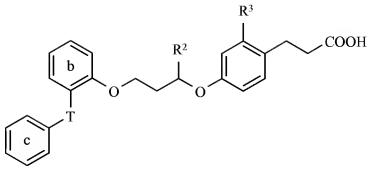


XIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein m is 1, 2, 3, or 4.

17. (Canceled).

18. (Previously Presented). The compound of Claim 16, wherein the compound is represented by structural formula XV,



XV

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

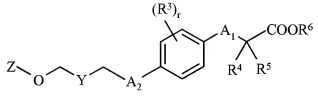
T is: a bond, O or C(O);

R^2 is: methyl, ethyl or cyclopropyl;

R^3 is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF_3, OCF_3, $N(CH_3)_2$, $S(O)_2CH_3$, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

19. (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula XVI,

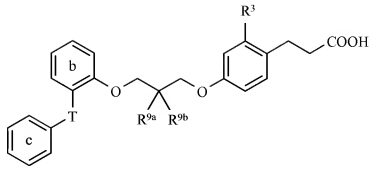


XVI

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein Y is a branched alkyl or C_3-C_6 cycloalkyl.

20. (Canceled).

21. (Previously Presented). The compound of Claim 19, wherein the compound structural formula XVIII,



XVIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

T is: a bond, O or C(O);

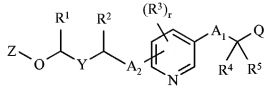
R^3 is: methyl or ethyl;

R^{9a} and R^{9b} are each independently hydrogen, methyl or ethyl, wherein at least one of R^{9a} and R^{9b} being methyl or ethyl;

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF_3, OCF_3, $S(O)_2CH_3$, $N(CH_3)_2$, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

22. (Canceled).

23. (Previously Presented). The compound of Claim 1, wherein the compound is a compound of formula XX,



XX

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 is: a bond, CH_2, O or S, and wherein A_1 and R^4 or A_1 and R^5 together being a 3- to 6-membered carbocyclyl when A_1 is a carbon;

A_2 is: O or S or CH_2;

Q is: $-C(O)OR^6$, or R^{6A};

Y is: a bond, C_1-C_6 alkyl or C_3-C_6 cycloalkyl;

Z is: a) aryl;

b) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,

c) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or

d) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and

wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R^7;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: 1, 2, 3, or 4;

R^1 and R^2 are each independently:

hydrogen,

haloalkyl,

C_1-C_6 alkyl,

$(CH_2)_nC_3$-C_8 cycloalkyl, or

R^1 and R^2 form a 4- to 8-membered nonaromatic carbocyclic ring; and wherein at least one of R^1 and R^2 is alkyl or cycloalkyl, and;

R^3 is: hydrogen,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

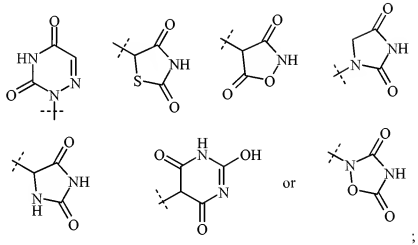
aryloxy,

C_1-C_6 alkyl,
C_1-C_6 alkoxy, or
C_3-C_8 cycloalkyl;

R^4 and R^5 are each independently: hydrogen or C_1-C_6 alkyl;

R^6 is: hydrogen, C_1-C_6 alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,



;

R^7 is: hydrogen,
oxo,
nitro,
cyano,
hydroxyl,
halo,
haloalkyl,
haloalkyloxy,
aryloxy,
arylalkyl,
aminoalkyl,
C_1-C_6 alkyl,
C_1-C_6 alkoxy,
$(CH_2)_nC_3$-C_8 cycloalkyl,
$C(O)R^9$,

$C(O)OR^9$,
$C(=NOR^8)R^9$,
$CR^8(OH)R^9$,
$C[=C(R^5)_2]R^9$,
OR^9,
SR^9 or
$S(O)_pR^9$;

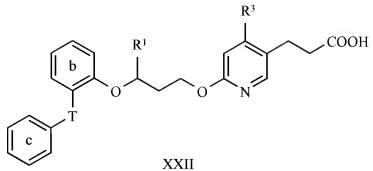
R^8 is: hydrogen or C_1-C_6 alkyl; and

R^9 is: hydrogen,
C_1-C_6 alkyl,
C_3-C_8 cycloalkyl,
aryl,
heteroaryl or
heterocyclyl,
wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:
hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy and C_3-C_8 cycloalkyl.

24. (Canceled).

25. (Canceled).

26. (Previously Presented). The compound of Claim 23, wherein the compound is a compound of structural formula XXII,



XXII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

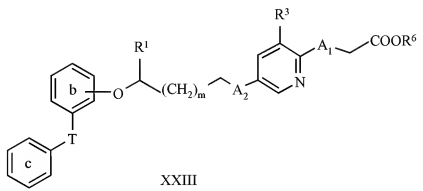
T is: a bond, -O- or -C(O)-;

R^1 is: methyl, ethyl or cyclopropyl;

R^3 is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF_3, OCF_3, $S(O)_2CH_3$, $N(CH_3)_2$, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

27. (Previously Presented). The compound of Claim 1, wherein the compound is a compound of structural formula XXIII,



XXIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 and A_2 are respectively:

O and O,

CH_2 and O,

CH_2 and S,

O and S or

S and O;

m is: 1, 2, 3 or 4;

R^1 is: C_1-C_3 alkyl; and

R^3 is: hydrogen, halo or C_1-C_6 alkyl;

R^6 and R^9 are each independently: hydrogen or C_1-C_6 alkyl;

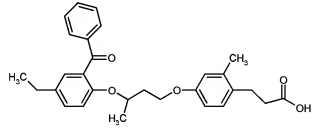
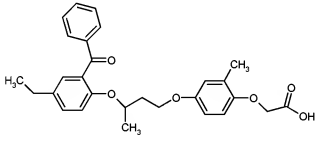
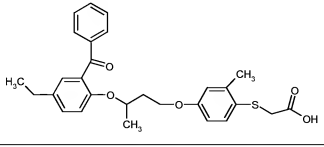
T is: a bond, -O-, -C(O)-, -S(O) –$S(O)_2$-, -C(=CH_2)-, -C(=NOH)- or -CH(OH)-; and

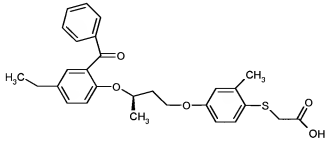
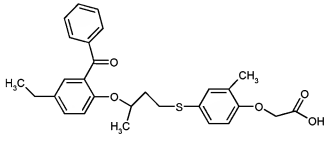
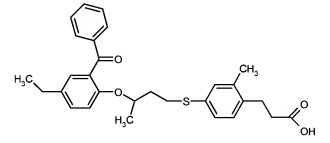
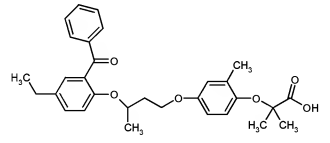
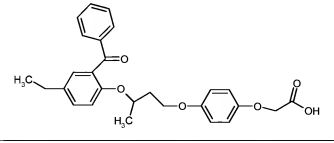
rings b and c are each optionally substituted with one or more groups independently selected from:

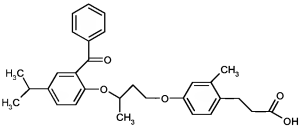
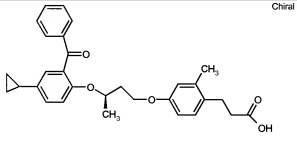
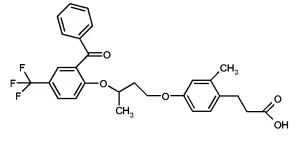
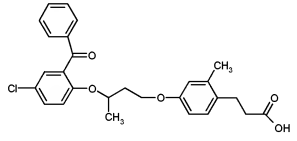
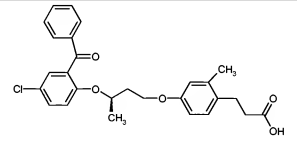
hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1-C_6 alkyl, C_1-C_6 alkoxy and $(CH_2)_nC_3$-C_8 cycloalkyl.

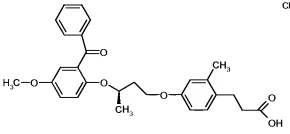
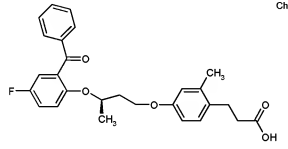
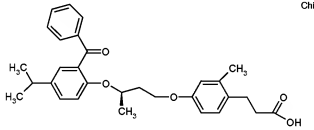
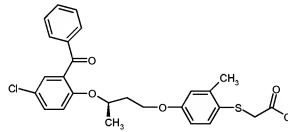
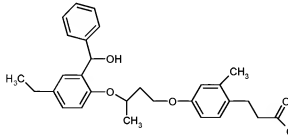
28. (Canceled).

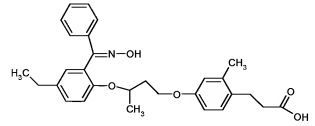
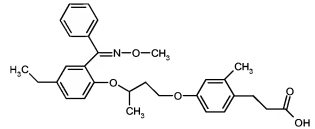
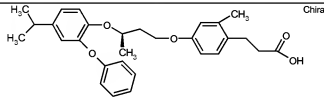
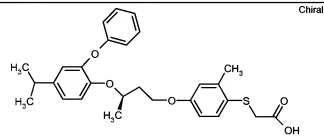
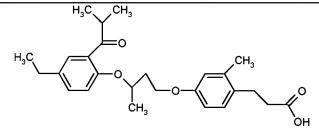
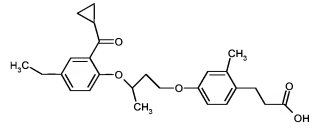
29. (Previously Presented). A compound of Claim 1 selected from the group consisting of:

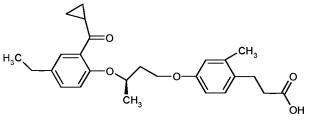
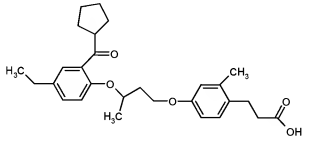
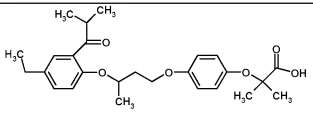
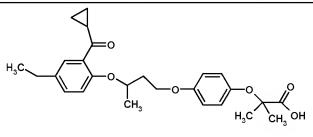
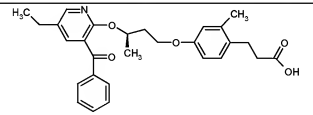
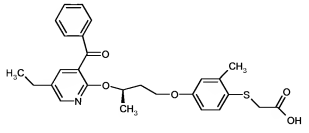
No.	Structure	Name
1		3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
2		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-acetic acid
3		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

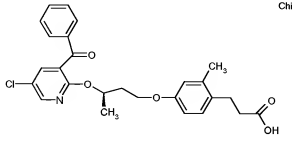
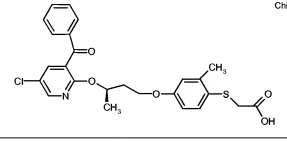
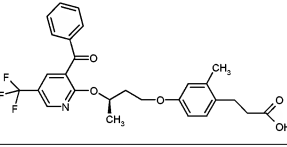
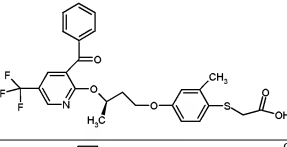
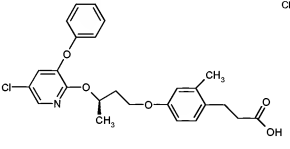
No.	Structure	Name
4		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
5		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenoxy}-acetic acid
6		3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic acid
7		2-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-2-methyl-propionic acid
8		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-acetic acid

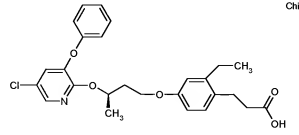
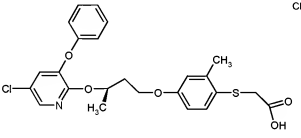
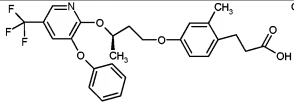
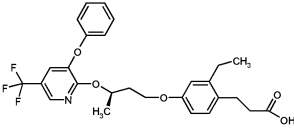
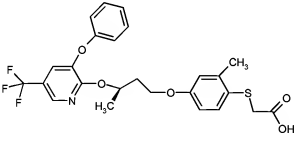
No.	Structure	Name
9		3-{4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
10	 Chiral	3-{4-[3-(2-Benzoyl-4-cyclopropyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
11		3-{4-[3-(2-Benzoyl-4-trifluoromethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
12		3-{4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
13		3-{4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

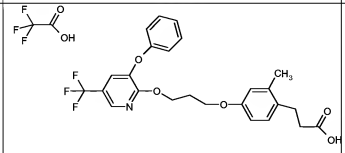
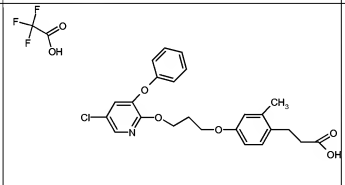
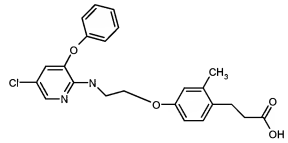
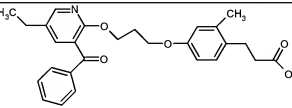
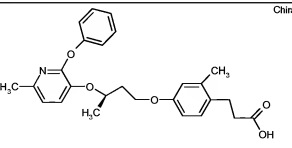
No.	Structure	Name
14	 Chiral	3-{4-[3-(2-Benzoyl-4-methoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
15	 Chiral	3-{4-[3-(2-Benzoyl-4-fluoro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
16	 Chiral	3-{4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
17		{4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
18		3-(4-{3-[4-Ethyl-2-(hydroxy-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid

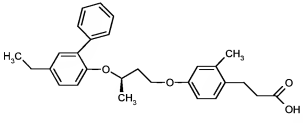
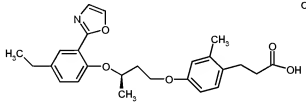
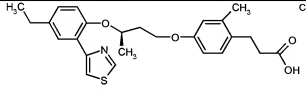
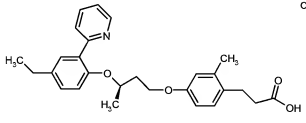
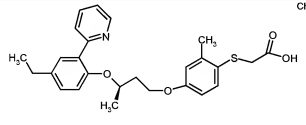
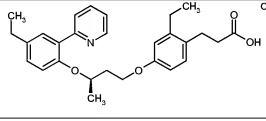
No.	Structure	Name
19		3-(4-{3-[4-Ethyl-2-(hydroxyimino-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
20		3-(4-{3-[4-Ethyl-2-(methoxyimino-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
21		3-{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
22		{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
23		3-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
24		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

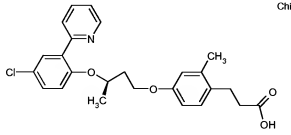
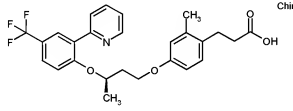
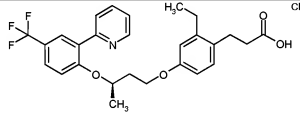
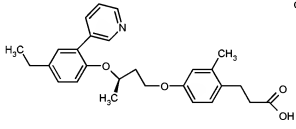
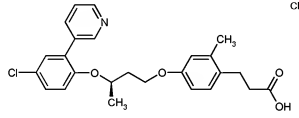
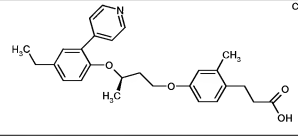
No.	Structure	Name
25		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
26		3-{4-[3-(2-Cyclopentanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
27		2-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
28		2-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
29		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
30		{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

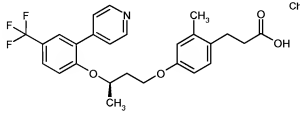
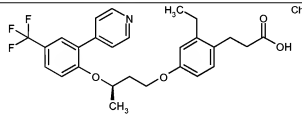
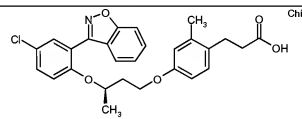
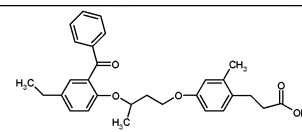
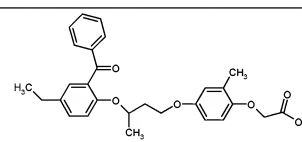
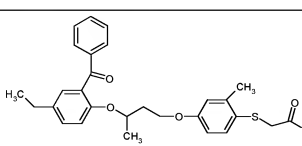
No.	Structure	Name
31	 Chiral	3-{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
32	 Chiral	{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
33	 Chiral	3-{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
34	 Chiral	{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
35	 Chiral	3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid

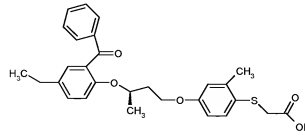
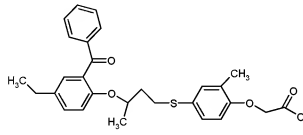
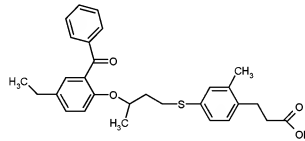
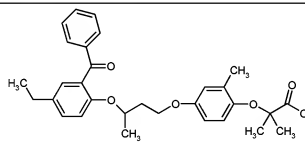
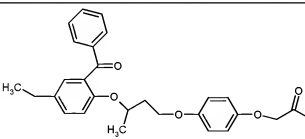
No.	Structure	Name
36	 Chiral	3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-ethyl-phenyl}-propionic acid
37	 Chiral	{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
38	 Chiral	3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
39	 Chiral	3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
40	 Chiral	3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid

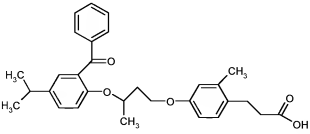
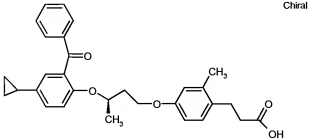
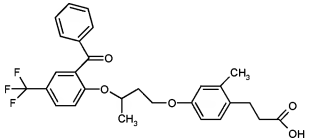
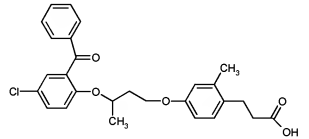
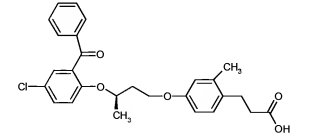
No.	Structure	Name
41		3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-propoxy]-phenyl}-propionic acid (trifluoroacetic acid salt)
42		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
43		3-{4-[2-(5-Chloro-3-phenoxy-pyridin-2-ylamino)-ethoxy]-2-methyl-phenyl}-propionic acid
44		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
45	Chiral 	3-{2-Methyl-4-[3-(6-methyl-2-phenoxy-pyridin-3-yloxy)-butoxy]-phenyl}-propionic acid

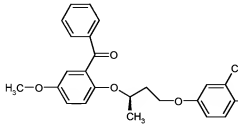
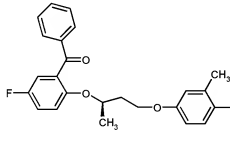
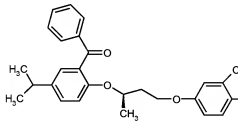
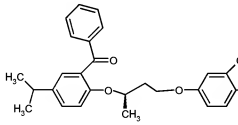
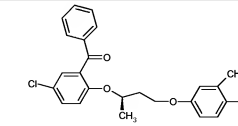
No.	Structure	Name
46		3-{4-[3-(5-Ethyl-biphenyl-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
47	Chiral 	3-{4-[3-(4-Ethyl-2-oxazol-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
48	Chiral 	3-{4-[3-(4-Ethyl-2-thiazol-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
49	Chiral 	3-{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
50	Chiral 	{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
51	Chiral 	3-{2-Ethyl-4-[3-(4-ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-phenyl}-propionic acid

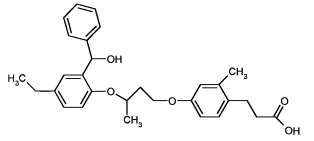
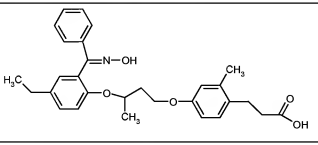
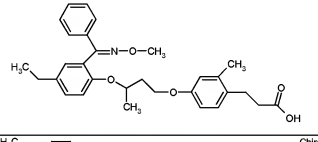
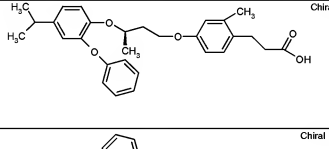
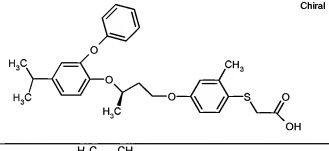
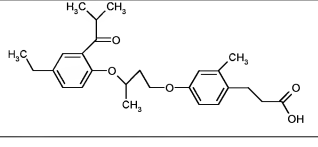
No.	Structure	Name
52	 Chiral	3-{4-[3-(4-Chloro-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
53	 Chiral	3-{2-Methyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
54	 Chiral	3-{2-Ethyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
55	 Chiral	3-{4-[3-(4-Ethyl-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
56	 Chiral	3-{4-[3-(4-Chloro-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
57	 Chiral	3-{4-[3-(4-Ethyl-2-pyridin-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

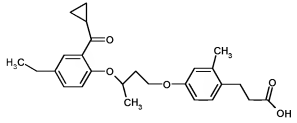
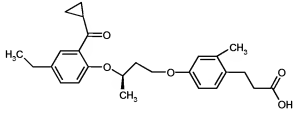
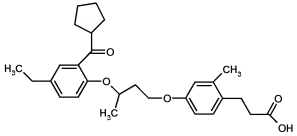
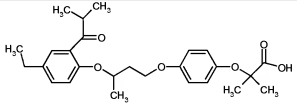
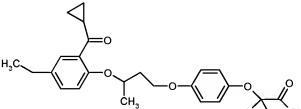
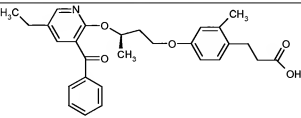
No.	Structure	Name
58		3-{2-Methyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
59		3-{2-Ethyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
60		3-{4-[3-(2-Benzo[d]isoxazol-3-yl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
61		3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
62		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-acetic acid
63		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

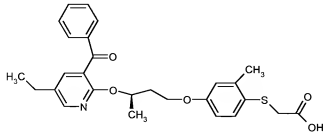
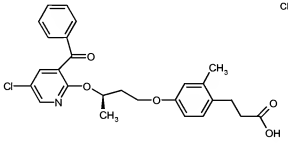
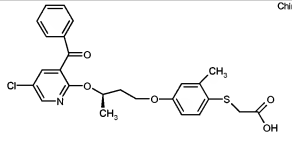
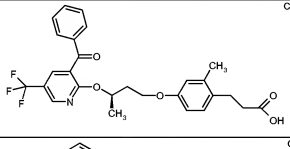
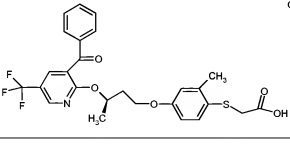
No.	Structure	Name
64		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
65		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenoxy}-acetic acid
66		3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic acid
67		2-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-2-methyl-propionic acid
68		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-acetic acid

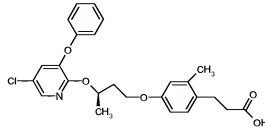
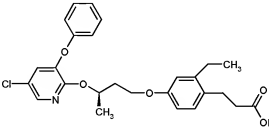
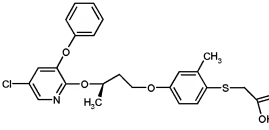
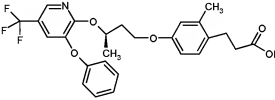
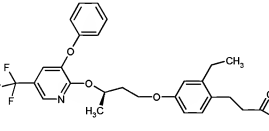
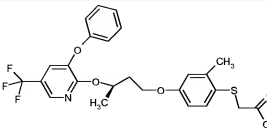
No.	Structure	Name
69		3-{4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
70	Chiral 	3-{4-[3-(2-Benzoyl-4-cyclopropyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
71		3-{4-[3-(2-Benzoyl-4-trifluoromethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
72		3-{4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
73		3-{4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

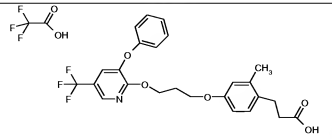
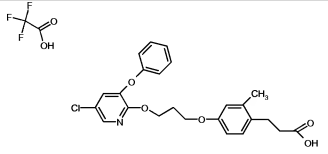
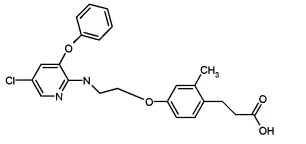
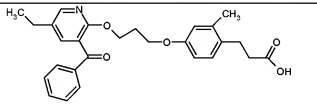
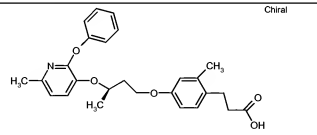
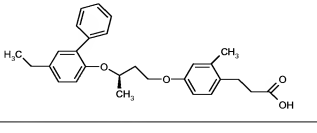
No.	Structure	Name
74	Chiral 	3-{4-[3-(2-Benzoyl-4-methoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
75	Chiral 	3-{4-[3-(2-Benzoyl-4-fluoro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
76	Chiral 	3-{4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
77	Chiral 	{4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
78		{4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

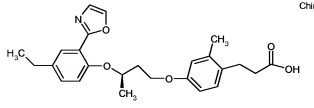
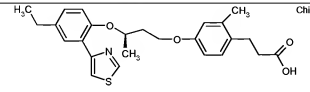
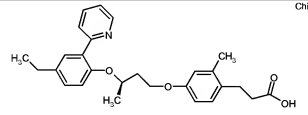
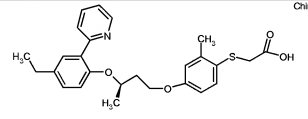
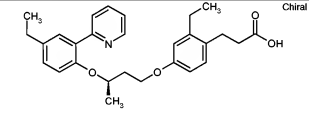
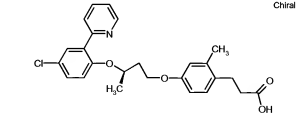
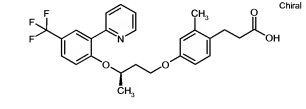
No.	Structure	Name
79		3-(4-{3-[4-Ethyl-2-(hydroxy-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
80		3-(4-{3-[4-Ethyl-2-(hydroxyimino-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
81		3-(4-{3-[4-Ethyl-2-(methoxyimino-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
82	 Chiral	3-{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
83	 Chiral	{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
84		3-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

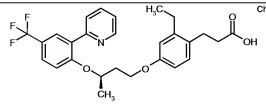
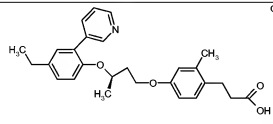
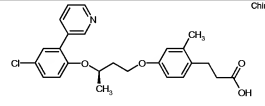
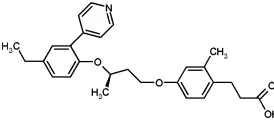
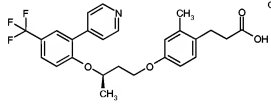
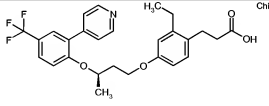
No.	Structure	Name
85		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
86		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
87		3-{4-[3-(2-Cyclopentanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
88		2-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
89		2-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
90		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid

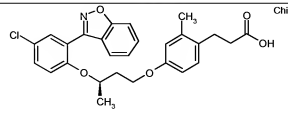
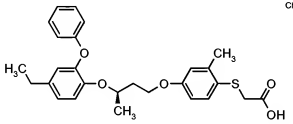
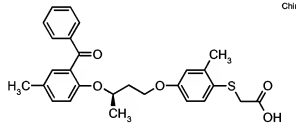
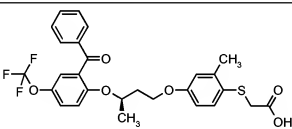
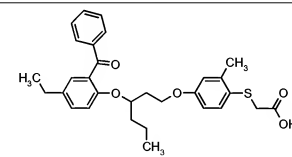
No.	Structure	Name
91		{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
92	 Chiral	3-{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
93	 Chiral	{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
94	 Chiral	3-{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
95	 Chiral	{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

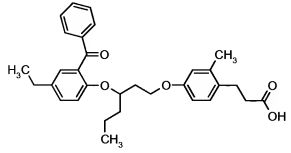
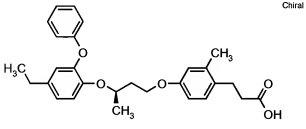
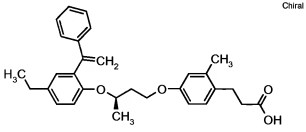
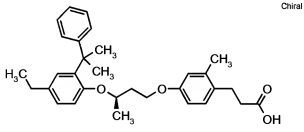
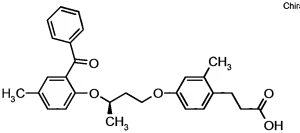
No.	Structure	Name
96	 Chiral	3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
97	 Chiral	3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-ethyl-phenyl}-propionic acid
98	 Chiral	{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
99	 Chiral	3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
100	 Chiral	3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
101	 Chiral	3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid

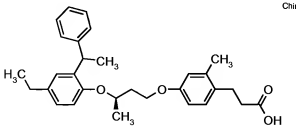
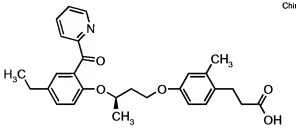
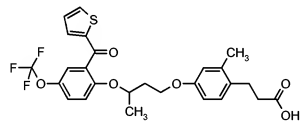
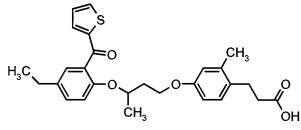
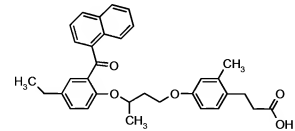
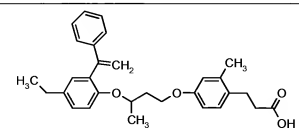
No.	Structure	Name
102		3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-propoxy]-phenyl}-propionic acid (trifluoroacetic acid salt)
103		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
104		3-{4-[2-(5-Chloro-3-phenoxy-pyridin-2-ylamino)-ethoxy]-2-methyl-phenyl}-propionic acid
105		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
106	 Chiral	3-{2-Methyl-4-[3-(6-methyl-2-phenoxy-pyridin-3-yloxy)-butoxy]-phenyl}-propionic acid
107		3-{4-[3-(5-Ethyl-biphenyl-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid

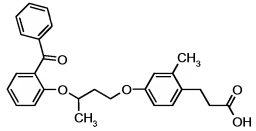
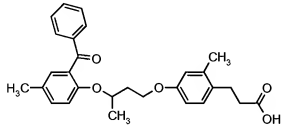
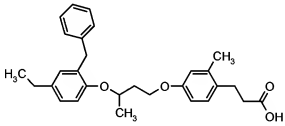
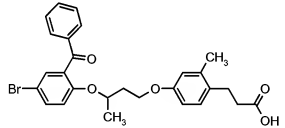
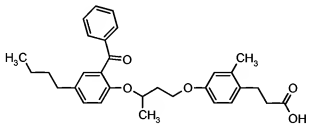
No.	Structure	Name
108	 Chiral	3-{4-[3-(4-Ethyl-2-oxazol-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
109	 Chiral	3-{4-[3-(4-Ethyl-2-thiazol-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
110	 Chiral	3-{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
111	 Chiral	{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
112	 Chiral	3-{2-Ethyl-4-[3-(4-ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-phenyl}-propionic acid
113	 Chiral	3-{4-[3-(4-Chloro-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
114	 Chiral	3-{2-Methyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-

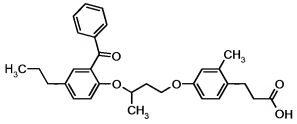
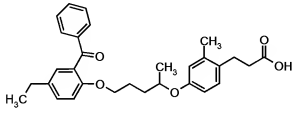
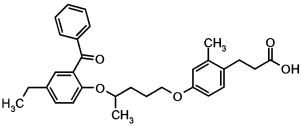
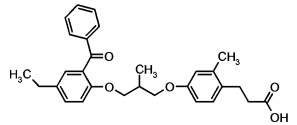
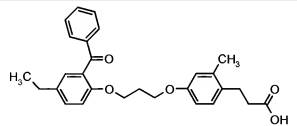
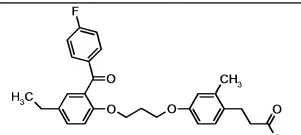
No.	Structure	Name
		phenyl}-propionic acid
115		3-{2-Ethyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
116		3-{4-[3-(4-Ethyl-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
117		3-{4-[3-(4-Chloro-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
118		3-{4-[3-(4-Ethyl-2-pyridin-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
119		3-{2-Methyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
120		3-{2-Ethyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid

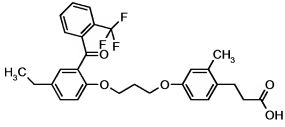
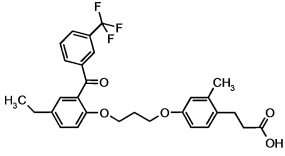
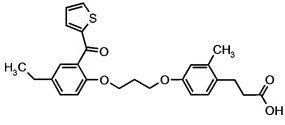
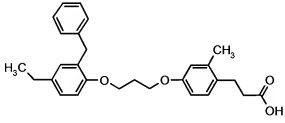
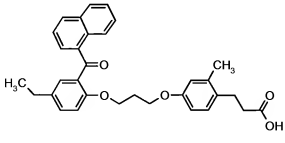
No.	Structure	Name
121	 Chiral	3-{4-[3-(2-Benzo[d]isoxazol-3-yl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
122	 Chiral	(R)-{4-[3-(4-ethyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
123	 Chiral	(R)-{4-[3-(2-benzoyl-4-methyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
124	 Chiral	(R)-{4-[3-(2-benzoyl-4-trifluoromethoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
125		{4-[3-(2-benzoyl-4-ethyl-phenoxy)-hexyloxy]-2-methyl-phenylsulfanyl}-acetic acid

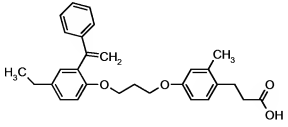
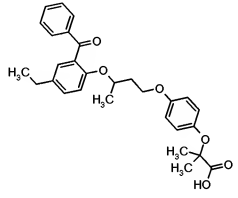
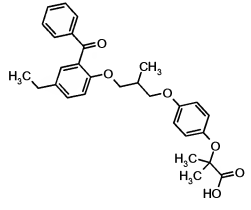
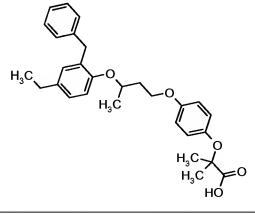
No.	Structure	Name
126		3-{4-[3-(2-benzoyl-4-ethyl-phenoxy)-hexyloxy]-2-methyl-phenyl}-propionic acid
127	 Chiral	(R)-3-{4-[3-(4-ethyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
128	 Chiral	(R)-3-(4-{3-[4-ethyl-2-(1-phenyl-vinyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
129	 Chiral	(R)-3-(4-{3-[4-ethyl-2-(1-methyl-1-phenyl-ethyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
130	 Chiral	(R)-3-{4-[3-(2-benzoyl-4-methyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

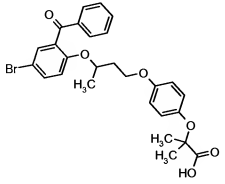
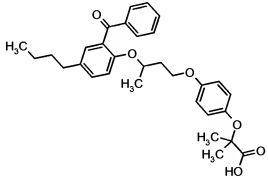
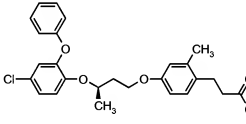
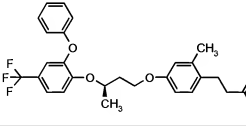
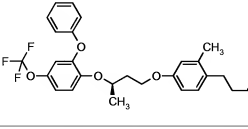
No.	Structure	Name
131	 Chiral	(R)-3-(4-{3-[4-ethyl-2-(1-phenyl-ethyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
132	 Chiral	(R)-3-(4-{3-[4-ethyl-2-(pyridine-2-carbonyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
133		3-(2-methyl-4-{3-[2-(thiophene-2-carbonyl)-4-trifluoromethoxy-phenoxy]-butoxy}-phenyl)-propionic acid
134		3-(4-{3-[4-ethyl-2-(thiophene-2-carbonyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
135		3-(4-{3-[4-ethyl-2-(naphthalene-1-carbonyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid
136		3-(4-{3-[4-ethyl-2-(1-phenyl-vinyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid

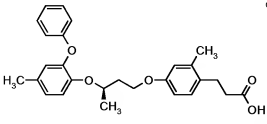
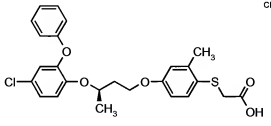
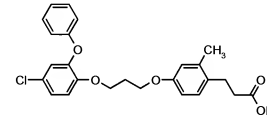
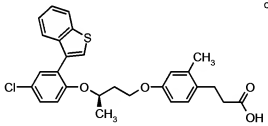
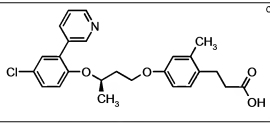
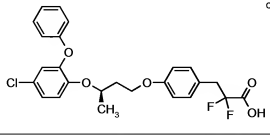
No.	Structure	Name
137		3-{4-[3-(2-benzoyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
138		3-{4-[3-(2-benzoyl-4-methyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
139		3-{4-[3-(2-benzyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
140		3-{4-[3-(2-benzoyl-4-bromo-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
141		3-{4-[3-(2-benzoyl-4-butyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

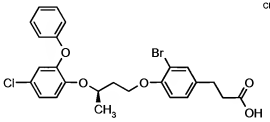
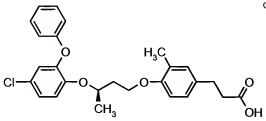
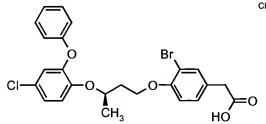
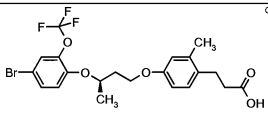
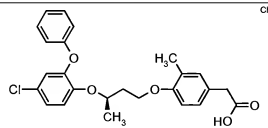
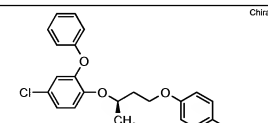
No.	Structure	Name
142		3-{4-[3-(2-benzoyl-4-propyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
143		3-{4-[4-(2-benzoyl-4-ethyl-phenoxy)-1-methyl-butoxy]-2-methyl-phenyl}-propionic acid
144		3-{4-[4-(2-benzoyl-4-ethyl-phenoxy)-pentyloxy]-2-methyl-phenyl}-propionic acid
145		3-{4-[3-(2-benzoyl-4-ethyl-phenoxy)-2-methyl-propoxy]-2-methyl-phenyl}-propionic acid
146		3-{4-[3-(2-benzoyl-4-ethyl-phenoxy)-propoxy]-2-methyl-phenyl}-propionic acid
147		3-(4-{3-[4-ethyl-2-(4-fluoro-benzoyl)-phenoxy]-propoxy}-2-methyl-phenyl)-propionic acid

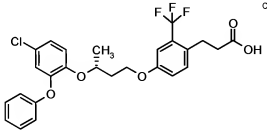
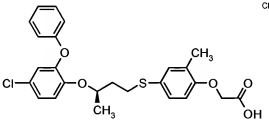
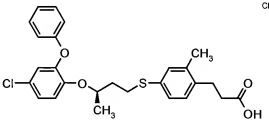
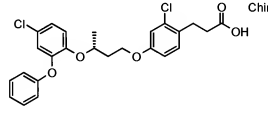
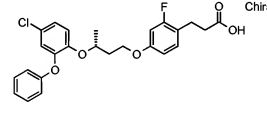
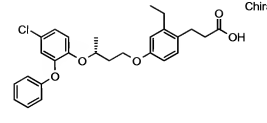
No.	Structure	Name
148		3-(4-{3-[4-ethyl-2-(2-trifluoromethyl-benzoyl)-phenoxy]-propoxy}-2-methyl-phenyl)-propionic acid
149		3-(4-{3-[4-ethyl-2-(3-trifluoromethyl-benzoyl)-phenoxy]-propoxy}-2-methyl-phenyl)-propionic acid
150		3-(4-{3-[4-ethyl-2-(thiophene-2-carbonyl)-phenoxy]-propoxy}-2-methyl-phenyl)-propionic acid
151		3-{4-[3-(2-benzyl-4-ethyl-phenoxy)-propoxy]-2-methyl-phenyl}-propionic acid
152		3-(4-{3-[4-ethyl-2-(naphthalene-1-carbonyl)-phenoxy]-propoxy}-2-methyl-phenyl)-propionic acid

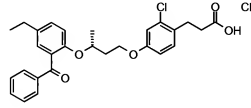
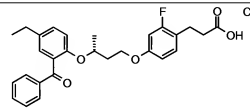
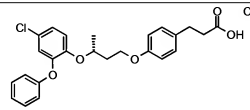
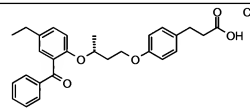
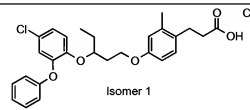
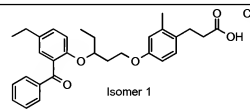
No.	Structure	Name
153		3-(4-{3-[4-ethyl-2-(1-phenyl-vinyl)-phenoxy]-propoxy}-2-methyl-phenyl)-propionic acid
154		2-{4-[3-(2-benzoyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
155		2-{4-[3-(2-benzoyl-4-ethyl-phenoxy)-2-methyl-propoxy]-phenoxy}-2-methyl-propionic acid
156		2-{4-[3-(2-benzyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid

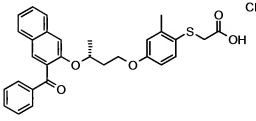
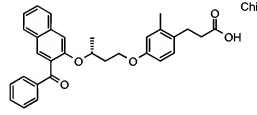
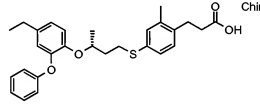
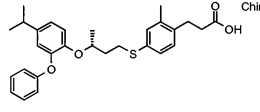
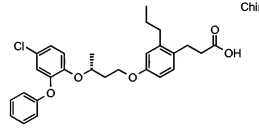
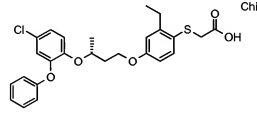
No.	Structure	Name
157		2-{4-[3-(2-benzoyl-4-bromo-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
158		2-{4-[3-(2-benzoyl-4-butyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
159	 Chiral	(R)- 3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
160	 Chiral	(R)-3-{2-methyl-4-[3-(2-phenoxy-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
161	 Chiral	(R)-3-{2-methyl-4-[3-(2-phenoxy-4-trifluoromethoxy-phenoxy)-butoxy]-phenyl}-propionic acid

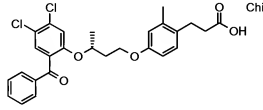
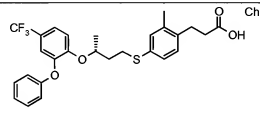
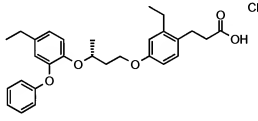
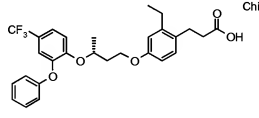
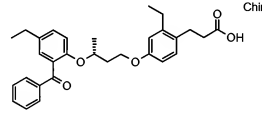
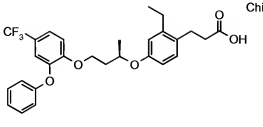
No.	Structure	Name
162		(R)-3-{2-methyl-4-[3-(4-methyl-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
163		(R)-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
164		3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-propoxy]-2-methyl-phenyl}-propionic acid
165		(R)-3-{4-[3-(2-benzo[b]thiophen-3-yl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
166		(R)- 3-{4-[3-(4-chloro-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
167		(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-2,2-difluoro-propionic acid

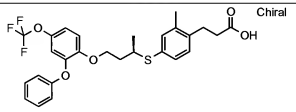
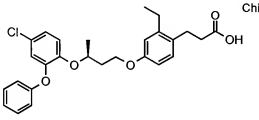
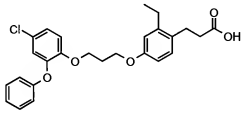
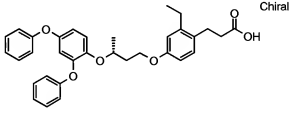
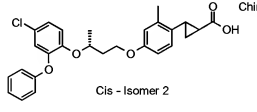
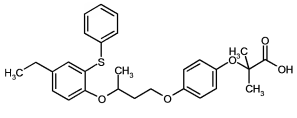
No.	Structure	Name
168		%)(R)-3-{3-bromo-4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
169		(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-3-methyl-phenyl}-propionic acid
170		(R)-{3-bromo-4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-acetic acid
171		(R)-3-{4-[3-(4-bromo-2-trifluoromethoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
172		(R)-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-3-methyl-phenyl}-acetic acid
173		(R)-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-acetic acid

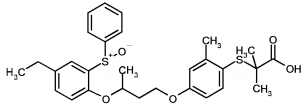
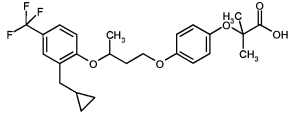
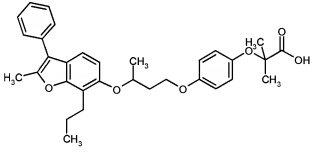
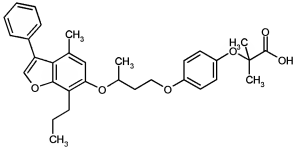
No.	Structure	Name
174	 Chiral	(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-2-trifluoromethyl-phenyl}-propionic acid
175	 Chiral	(R)-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenoxy}-acetic acid
176	 Chiral	(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic acid
177	 Chiral	(R)-3-{2-Chloro-4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
178	 Chiral	(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-fluoro-phenyl}-propionic acid
179	 Chiral	(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic acid

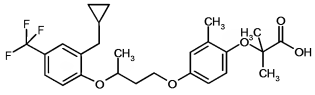
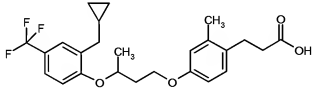
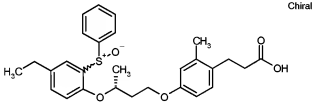
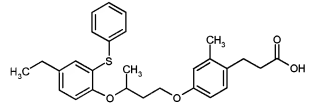
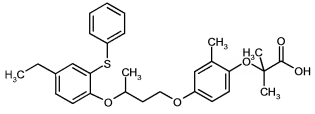
No.	Structure	Name
180		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-chloro-phenyl}-propionic acid
181		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-fluoro-phenyl}-propionic acid
182		(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
183		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-phenyl}-propionic acid
184	 Isomer 1	(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-pentyloxy]-2-methyl-phenyl}-propionic acid
185	 Isomer 1	(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-pentyloxy]-2-methyl-phenyl}-propionic acid

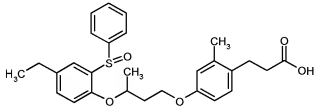
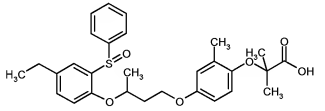
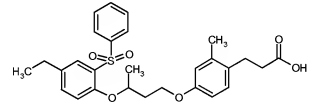
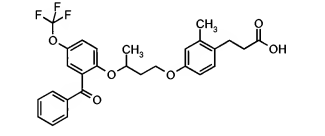
No.	Structure	Name
186	 Chiral	(R)-{4-[3-(3-Benzoyl-naphthalen-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
187	 Chiral	(R)-3-{4-[3-(3-Benzoyl-naphthalen-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
188	 Chiral	(R)-3-{4-[3-(4-Ethyl-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic acid
189	 Chiral	(R)-3-{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic acid
190	 Chiral	(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-propyl-phenyl}-propionic acid
191	 Chiral	(R)-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-ethyl-phenylsulfanyl}-acetic acid

No.	Structure	Name
192		(R)-3-{4-[3-(2-Benzoyl-4,5-dichloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
193		(R)-3-{2-Methyl-4-[3-(2-phenoxy-4-trifluoromethyl-phenoxy)-butylsulfanyl]-phenyl}-propionic acid
194		(R)-3-{2-Ethyl-4-[3-(4-ethyl-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
195		(R)-3-{2-Ethyl-4-[3-(2-phenoxy-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
196		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic acid
197		(R)-3-{2-Ethyl-4-[1-methyl-3-(2-phenoxy-4-trifluoromethyl-phenoxy)-propoxy]-phenyl}-propionic acid

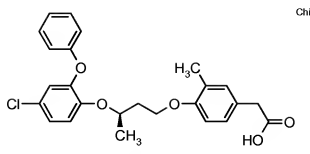
No.	Structure	Name
198	 Chiral	(R)-3-{2-Methyl-4-[1-methyl-3-(2-phenoxy-4-trifluoromethoxy-phenoxy)-propylsulfanyl]-phenyl}-propionic acid
199	 Chiral	(S)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic acid
200		3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-propoxy]-2-ethyl-phenyl}-propionic acid
201	 Chiral	(R)-3-{4-[3-(2,4-Diphenoxy-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic acid
202	 Chiral Cis - Isomer 2	2-{4-[4-(4-Chloro-2-phenoxy-phenyl)-3-methyl-butoxy]-2-methyl-phenyl}-cyclopropanecarboxylic acid
203		(R, S)-2-{4-[3-(4-Ethyl-2-phenylsulfanyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid

No.	Structure	Name
204		2-{4-[3-(R,S-2-Benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-2-methyl-propionic acid (enamtiomer pair 1)
205		(R, S)-2-{4-[3-(2-Cyclopropylmethyl-4-trifluoromethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
206		(R, S)-2-Methyl-2-{4-[3-(2-methyl-3-phenyl-7-propyl-benzofuran-6-yloxy)-butoxy]-phenoxy}-propionic acid
207		(R, S)-2-Methyl-2-{4-[3-(4-methyl-3-phenyl-7-propyl-benzofuran-6-yloxy)-butoxy]-phenoxy}-propionic acid

No.	Structure	Name
208		(R, S)-2-{4-[3-(2-Cyclopropylmethyl-4-trifluoromethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-2-methyl-propionic acid
209		(R, S)-3-{4-[3-(2-Cyclopropylmethyl-4-trifluoromethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
210	 Chiral	3-{R-4-[3-(R, S-2-Benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
211		3-{4-[3-(4-Ethyl-2-phenylsulfanyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid isomer 2
212		(R, S)-2-{4-[3-(4-Ethyl-2-phenylsulfanyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid

No.	Structure	Name
213		(R, S)-3-{4-[3-(R, S-2-Benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
214		(R, S)-2-{4-[3-(R, S-2-Benzenesulfinyl-4-ethyl-phenoxy) -butoxy]-2-methyl-phenoxy}-2-methyl-propionic acid
215		(R, S)-3-{4-[3-(2-Benzenesulfonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
216		3-{4-[3-(2-Benzoyl-4-trifluoromethoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

30. (Original). The compound of Claim 29, wherein the compound is



or a pharmaceutically acceptable salt, solvate or hydrate thereof.

31. (Previously Presented). A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a compound of Claim 1 or a pharmaceutically acceptable salt, solvate or hydrate thereof.

32. (Canceled).

33. (Canceled).

34. (Canceled).

35. (Canceled).

36. (Canceled).

37. (Canceled).

38. (Canceled).

39. (Canceled).

40. (Canceled).

41. (Canceled).

42. (Canceled).

43. (Previously Presented). A method for lowering blood-glucose in a mammal in need thereof comprising the step of administering an effective amount of a compound of Claim 1.

44. (Canceled).

45. (Canceled).

46. (Canceled).

47. (Canceled).

48. (Canceled).

49. 49. (Canceled)